LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030623\
Data File : BG056851.D
Acq On : 6 Mar 2023 21:33
Operator : CG/JU
Sample : 01712-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZX8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Title : SVOA CALIBRATION

Signal : TIC: BG056851.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.254	118	122	127	rVB2	8606	12257	3.64%	0.286%
2	5.435	321	323	329	rVB3	6884	10757	3.20%	0.251%
3	7.538	675	681	692	rVB2	45608	82475	24.51%	1.924%
4	7.720	706	712	721	rBV	32160	54629	16.23%	1.275%
5	7.938	744	749	757	rVB	42924	73286	21.78%	1.710%
6	8.414	824	830	840	rVB	70742	121633	36.14%	2.838%
7	9.101	942	947	954	rBV2	37570	65902	19.58%	1.538%
8	9.242	966	971	977	rBV2	14143	24990	7.43%	0.583%
9	9.589	1024	1030	1037	rBV2	51457	84803	25.20%	1.979%
10	9.971	1092	1095	1098	rVB3	1213	1323	0.39%	0.031%
11	10.317	1148	1154	1163	rVB	41161	72187	21.45%	1.684%
12	10.382	1163	1165	1167	rBV	685	576	0.17%	0.013%
13	10.453	1170	1177	1183	rBV3	1809	4202	1.25%	0.098%
14	10.605	1201	1203	1206	rBV	664	710	0.21%	0.017%
15	10.752	1224	1228	1231	rBV2	827	938	0.28%	0.022%
16	10.858	1240	1246	1253	rVB2	58009	99886	29.68%	2.331%
17	11.163	1296	1298	1301	rBV2	538	590	0.18%	0.014%
18	11.257	1306	1314	1322	rVB	105555	190101	56.49%	4.436%
19	11.375	1329	1334	1346	rVB2	17832	31313	9.30%	0.731%
20	11.757	1398	1399	1405	rVB3	843	1110	0.33%	0.026%
21	12.010	1440	1442	1448	rVB2	1282	2048	0.61%	0.048%
22	12.315	1491	1494	1497	rBV2	391	554	0.16%	0.013%
23	12.474	1519	1521	1525	rBV2	431	632	0.19%	0.015%
24	12.609	1540	1544	1547	rBV3	1235	1440	0.43%	0.034%
25	12.809	1574	1578	1580	rBV2	1074	1237	0.37%	0.029%
26	13.108	1626	1629	1632	rVB2	411	603	0.18%	0.014%
27	13.155	1632	1637	1638	rBV2	523	554	0.16%	0.013%
28	13.214	1643	1647	1650	rBV	308	607	0.18%	0.014%
29	13.819	1745	1750	1753	rBV4	1695	2504	0.74%	0.058%
30	13.878	1756	1760	1761	rVB2	517	516	0.15%	0.012%
31	13.943	1769	1771	1776	rVB2	476	600	0.18%	0.014%
32	14.007	1778	1782	1784	rBV	467	510	0.15%	0.012%
33	14.407	1844	1850	1856	rBV	110439	158560	47.12%	3.700%
34	14.718	1896	1903	1910	rBV	122043	182986	54.38%	4.270%
35	14.871	1925	1929	1935	rVB	14947	19236	5.72%	0.449%
36	15.024	1948	1955	1960	rVB	168826	269782	80.17%	6.295%

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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DBZX8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Title : SVOA CALIBRATION

37	15.176	1975	1981	1988	rBV2	42532	67242	19.98%	1.569%
38	15.335	2004	2008	2012	rVB4	2046	3344	0.99%	0.078%
39	15.558	2045	2046	2049	rVB	815	555	0.16%	0.013%
40	15.717	2070	2073	2076	rBV	533	653	0.19%	0.015%
41	15.811	2087	2089	2093	rVB2	1330	1633	0.49%	0.038%
42	15.999	2114	2121	2128	rVV	155706	235417	69.96%	5.493%
43	16.105	2133	2139	2148	rVB	34917	52026	15.46%	1.214%
44	16.352	2175	2181	2187	rVB	42346	58667	17.43%	1.369%
45	16.422	2189	2193	2195	rBV2	489	513	0.15%	0.012%
46	16.510	2203	2208	2211	rBV3	4320	5573	1.66%	0.130%
47	16.545	2211	2214	2220	rVB5	5166	8486	2.52%	0.198%
48	16.675	2231	2236	2240	rBV	10503	12510	3.72%	0.292%
49	16.739	2246	2247	2251	rVB2	738	615	0.18%	0.014%
50	16.798	2253	2257	2259	rVB	642	844	0.25%	0.020%
51	16.939	2278	2281	2286	rVB2	726	840	0.25%	0.020%
52	17.080	2302	2305	2310	rVB2	9149	12090	3.59%	0.282%
53	17.274	2334	2338	2343	rVB2	924	1541	0.46%	0.036%
54	17.339	2343	2349	2352	rBV	5762	7323	2.18%	0.171%
55	17.468	2368	2371	2374	rVB	935	948	0.28%	0.022%
56	17.550	2383	2385	2388	rVB	904	922	0.27%	0.022%
57	17.685	2406	2408	2412	rVB	757	560	0.17%	0.013%
58	17.756	2413	2420	2426	rVV	214299	317998	94.50%	7.420%
59	17.850	2429	2436	2446	rBV2	153352	231024	68.65%	5.391%
60	17.950	2451	2453	2456	rBV	599	705	0.21%	0.016%
61	17.979	2456	2458	2461	rVB	707	782	0.23%	0.018%
62	18.038	2465	2468	2471	rVB2	1948	2096	0.62%	0.049%
63	18.126	2481	2483	2488	rVB2	733	1033	0.31%	0.024%
64	18.191	2488	2494	2497	rVB2	1011	1289	0.38%	0.030%
65	18.238	2499	2502	2504	rBV	859	825	0.25%	0.019%
66	18.332	2516	2518	2519	rBV	791	515	0.15%	0.012%
67	18.384	2525	2527	2530	rVB	661	626	0.19%	0.015%
68	18.478	2540	2543	2544	rBV	788	665	0.20%	0.016%
69	18.490	2544	2545	2549	rVV	1298	1020	0.30%	0.024%
70	18.590	2558	2562	2568	rBV2	41040	59520	17.69%	1.389%
71	18.731	2584	2586	2590	rBV3	887	1243	0.37%	0.029%
72	18.807	2597	2599	2601	rVB2	944	685	0.20%	0.016%
73	18.872	2608	2610	2613	rBV2	902	840	0.25%	0.020%
74	18.960	2623	2625	2626	rBV2	1315	1053	0.31%	0.025%
75	19.007	2630	2633	2636	rBV2	5034	5074	1.51%	0.118%
76	19.054	2638	2641	2642	rBV2	646	701	0.21%	0.016%

LSC Area Percent Report

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Sample : 01712-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBZX8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Title : SVOA CALIBRATION

77	19.107	2648	2650	2651	rBV2	1384	1313	0.39%	0.031%
78	19.178	2659	2662	2669	rVB4	3751	4857	1.44%	0.113%
79	19.395	2695	2699	2703	rBV4	4083	5220	1.55%	0.122%
80	19.442	2705	2707	2710	rVB4	2183	2686	0.80%	0.063%
81	19.471	2710	2712	2713	rBV	1063	979	0.29%	0.023%
82	19.571	2727	2729	2730	rBV	1966	1113	0.33%	0.026%
83	19.595	2731	2733	2734	rBV	2277	1505	0.45%	0.035%
84	19.842	2772	2775	2785	rBV5	12290	17063	5.07%	0.398%
85	19.953	2792	2794	2797	rBV4	2297	1901	0.56%	0.044%
86	19.983	2797	2799	2802	rBV3	2197	2456	0.73%	0.057%
87	20.112	2815	2821	2828	rVB	86298	124271	36.93%	2.900%
88	20.535	2889	2893	2896	rBV2	9672	12700	3.77%	0.296%
89	20.588	2900	2902	2906	rVB3	6976	7360	2.19%	0.172%
90	21.023	2974	2976	2978	rBV3	4245	4766	1.42%	0.111%
91	21.058	2980	2982	2985	rVB	9648	12300	3.66%	0.287%
92	21.645	3077	3082	3090	rVB	125376	199998	59.43%	4.667%
93	22.062	3146	3153	3165	rVB	198030	336521	100.00%	7.852%
94	22.280	3185	3190	3198	rVV	78382	129144	38.38%	3.013%
95	22.356	3198	3203	3209	rVV	117779	190043	56.47%	4.434%
96	22.421	3209	3214	3217	rVV2	37898	62819	18.67%	1.466%
97	22.456	3218	3220	3232	rVB4	42800	74868	22.25%	1.747%
98	23.108	3327	3331	3336	rVB7	13631	25701	7.64%	0.600%
99	25.335	3705	3710	3720	rVB4	19222	54947	16.33%	1.282%
100	25.588	3743	3753	3765	rVB	108661	335022	99.55%	7.817%

Sum of corrected areas: 4285586

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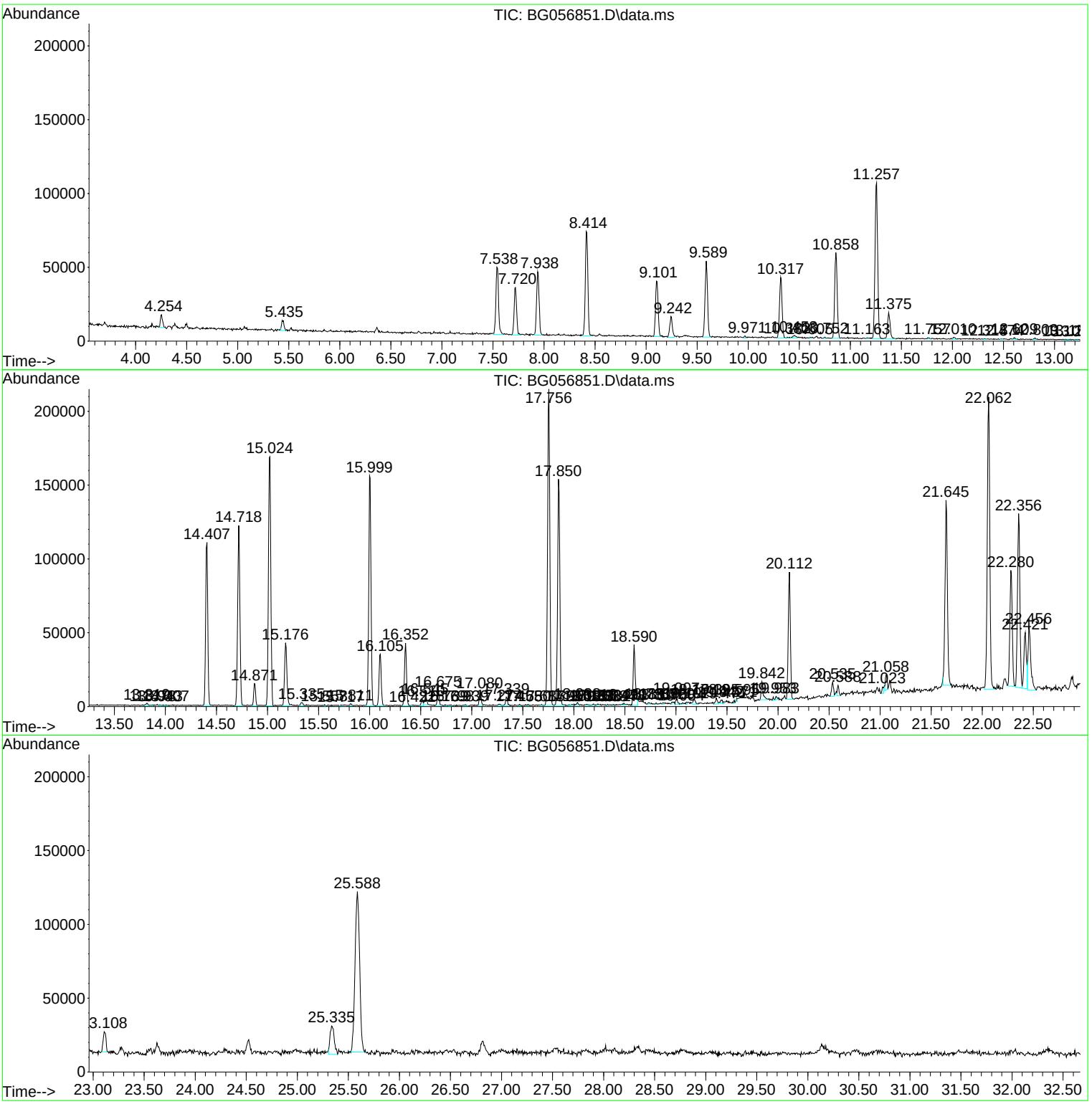
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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Quant Title : SVOA CALIBRATION

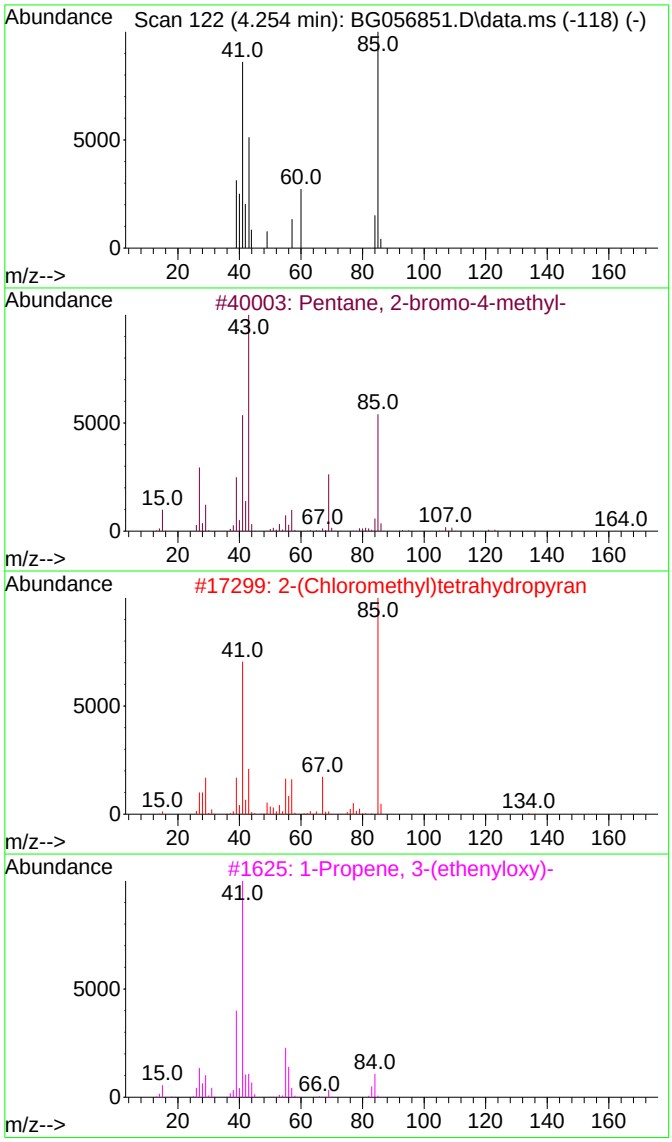
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

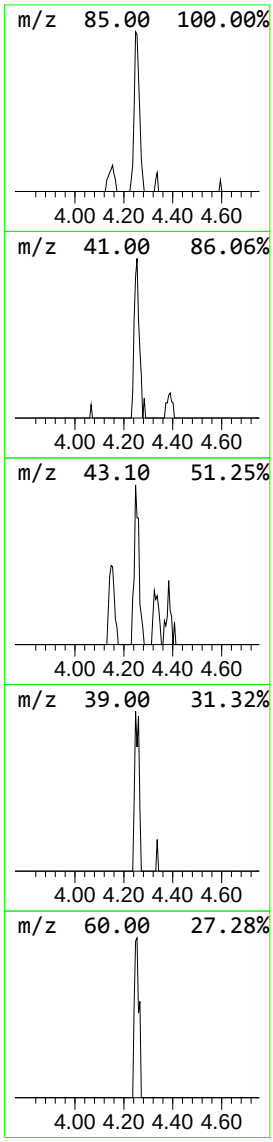
****

Peak Number 1 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.254	2.02 ng/ul	12257	1,4-Dichlorobenzene-d4	8.420

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-bromo-4-methyl-	164	C6H13Br	030310-22-6	39
2	2-(Chloromethyl)tetrahydropyran	134	C6H11ClO	018420-41-2	36
3	1-Propene, 3-(ethenyloxy)-	84	C5H8O	003917-15-5	9
4	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9
5	2-Pyrrolidinone	85	C4H7NO	000616-45-5	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030623\
Data File : BG056851.D
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Instrument :
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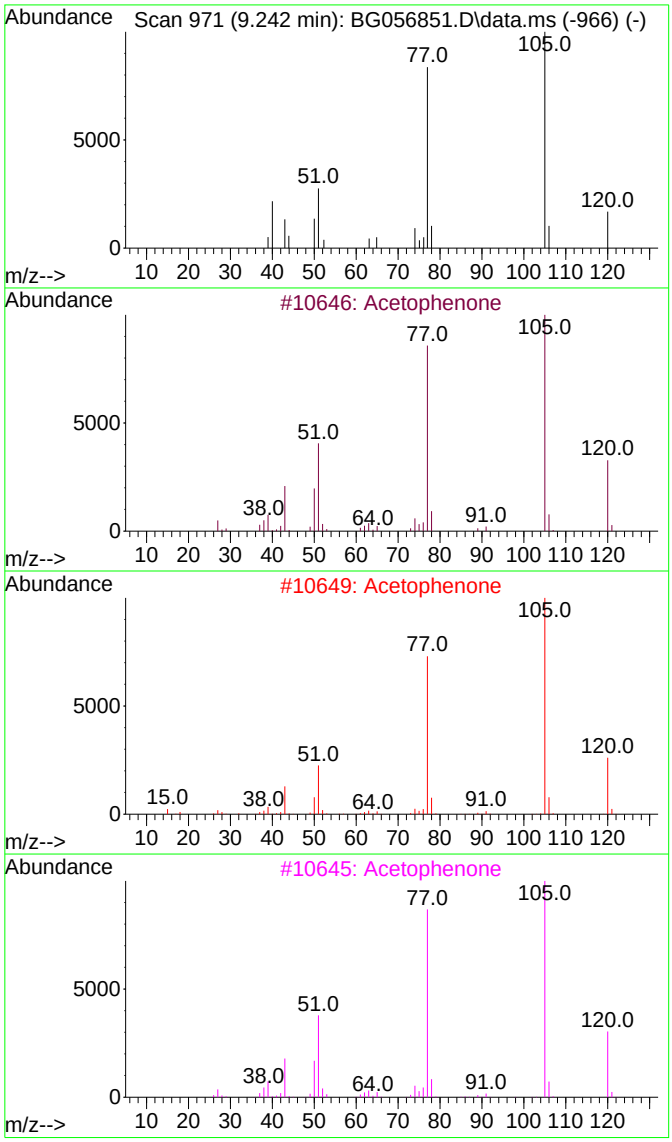
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

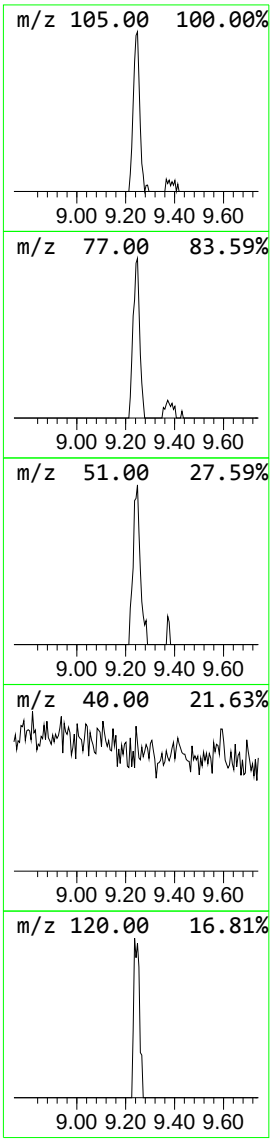
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Acetophe none Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.242	4.11 ng/ul	24990	1,4-Dichlorobenzene-d4	8.420

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone	120	C8H8O	000098-86-2	91
2	Acetophenone	120	C8H8O	000098-86-2	86
3	Acetophenone	120	C8H8O	000098-86-2	83
4	Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	83
5	N-cyanobenzamide	146	C8H6N2O	015150-25-1	83





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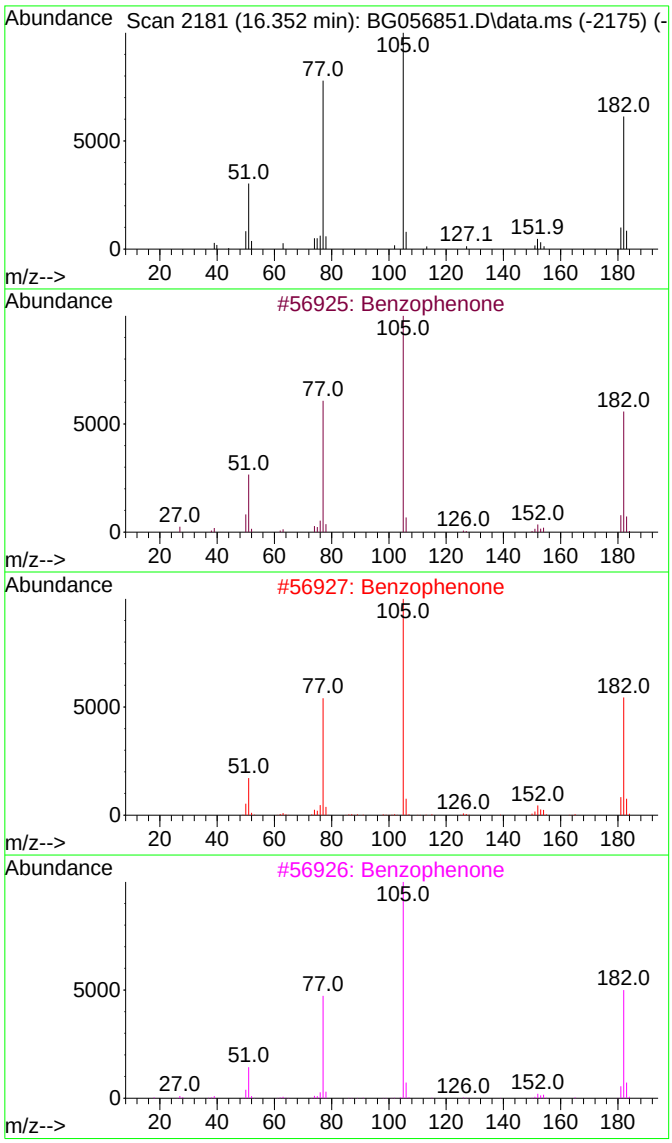
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

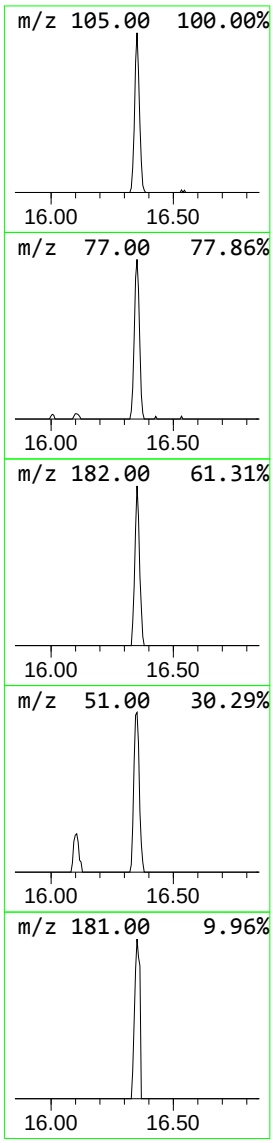
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.352	4.35 ng/ul	58667	Acenaphthene-d10	15.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Benzophenone	182	C13H10O	000119-61-9	93
3	Benzophenone	182	C13H10O	000119-61-9	91
4	Benzophenone	182	C13H10O	000119-61-9	91
5	Benzophenone	182	C13H10O	000119-61-9	91





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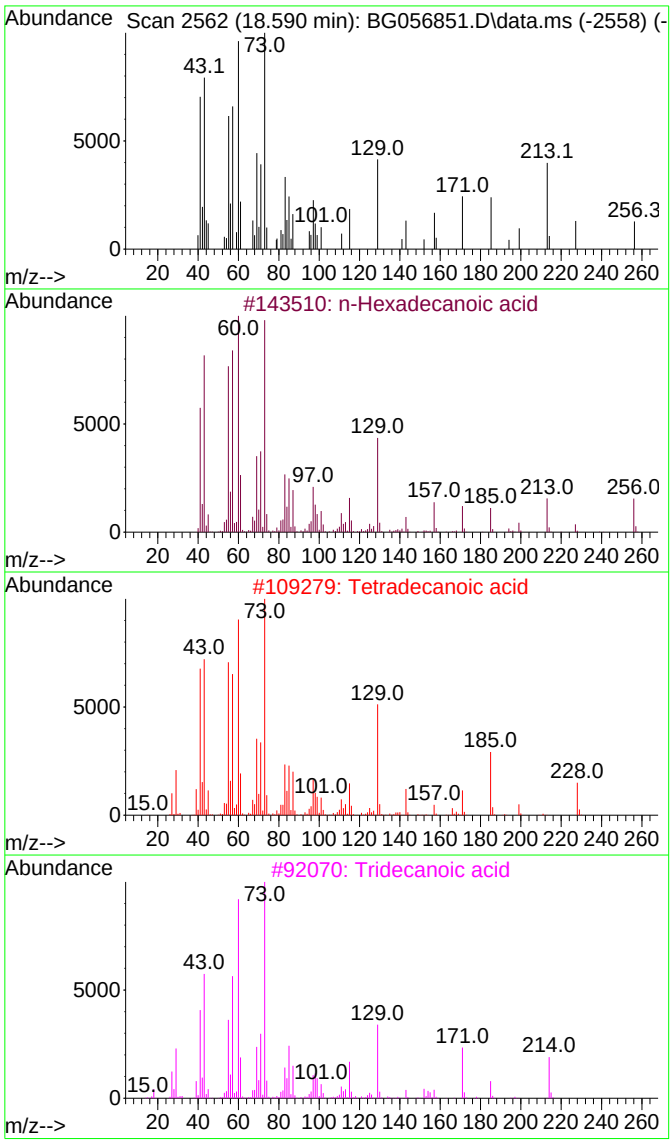
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

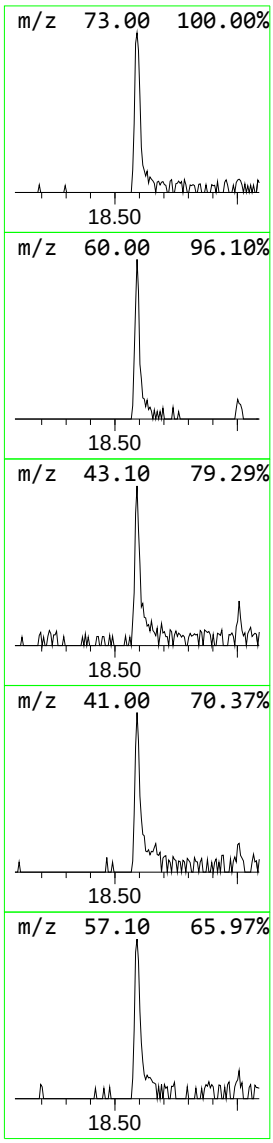
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.590	3.74 ng/ul	59520	Phenanthrene-d10	17.756

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
3	Tridecanoic acid	214	C13H26O2	000638-53-9	72
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	64





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ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZX8

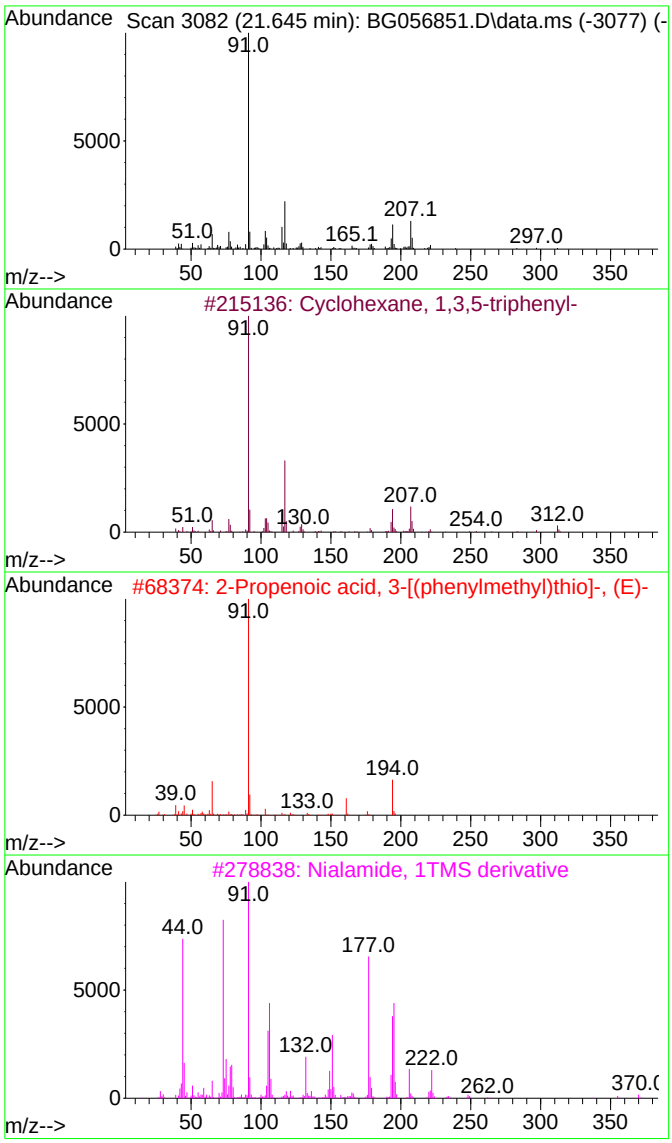
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

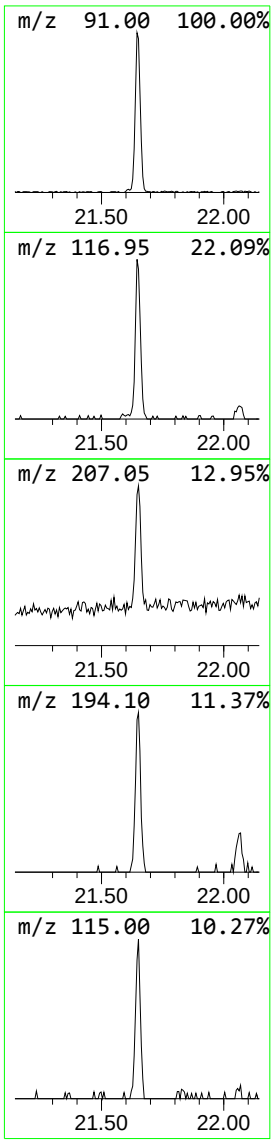
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Cyclohexane, 1,3,5-triphenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.645	11.89 ng/ul	199998	Chrysene-d12	22.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3,5-triphenyl-	312	C24H24	028336-57-4	86
2	2-Propenoic acid, 3-[(phenylmeth...	194	C10H10O2S	013831-01-1	22
3	Nialamide, 1TMS derivative	370	C19H26N4O2Si	1000485-79-3	22
4	Carbonic acid, monoamide, N-benz...	207	C12H17NO2	1000451-48-5	12
5	N-[4-(Benzyloxy)-2-nitrophenyl]a...	286	C15H14N2O4	026697-34-7	10





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Data File : BG056851.D
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZX8

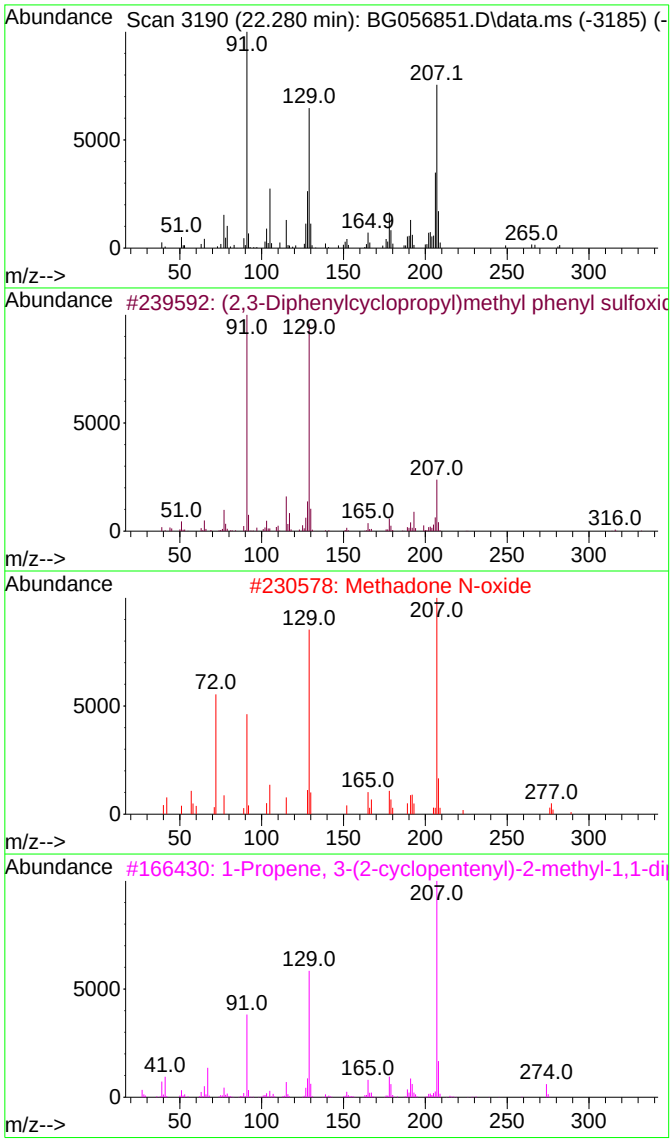
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

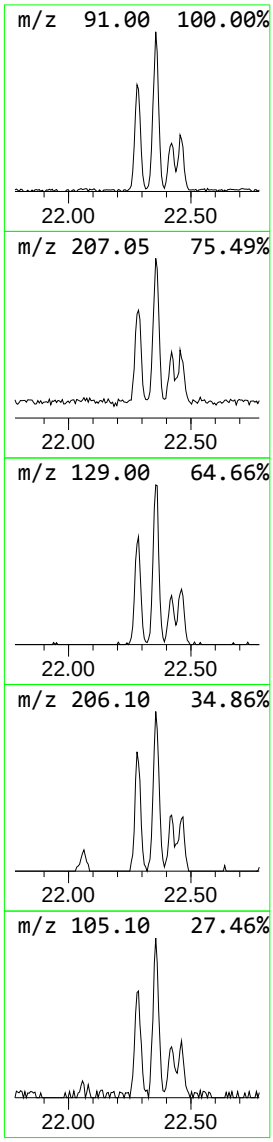
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (2,3-Diphenylcyclopropyl)me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.280	7.68 ng/ul	129144	Chrysene-d12	22.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	78
2	Methadone N-oxide	325	C21H27NO2	1000120-80-7	59
3	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1010154-23-3	59
4	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	53
5	1H-Indole, 5-methyl-2-phenyl-	207	C15H13N	013228-36-9	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030623\
Data File : BG056851.D
Acq On : 6 Mar 2023 21:33
Operator : CG/JU
Sample : 01712-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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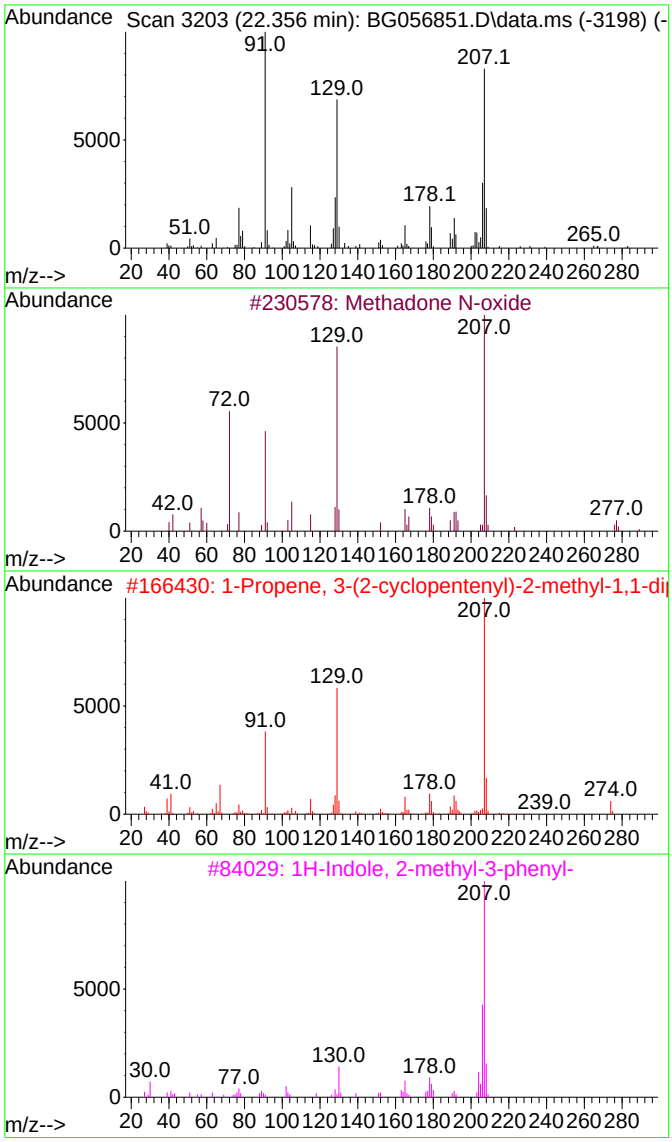
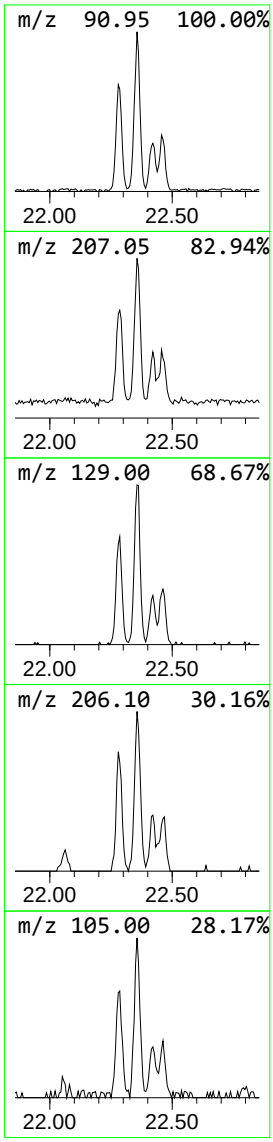
Peak Number 7 Methadone N-oxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.356	11.29 ng/ul	190043	Chrysene-d12	22.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methadone N-oxide	325	C21H27NO2	1000120-80-7	59
2	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1010154-23-3	59
3	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	49
4	2-(4-Methylphenyl)-1H-indole	207	C15H13N	055577-25-8	49
5	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030623\
Data File : BG056851.D
Acq On : 6 Mar 2023 21:33
Operator : CG/JU
Sample : 01712-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZX8

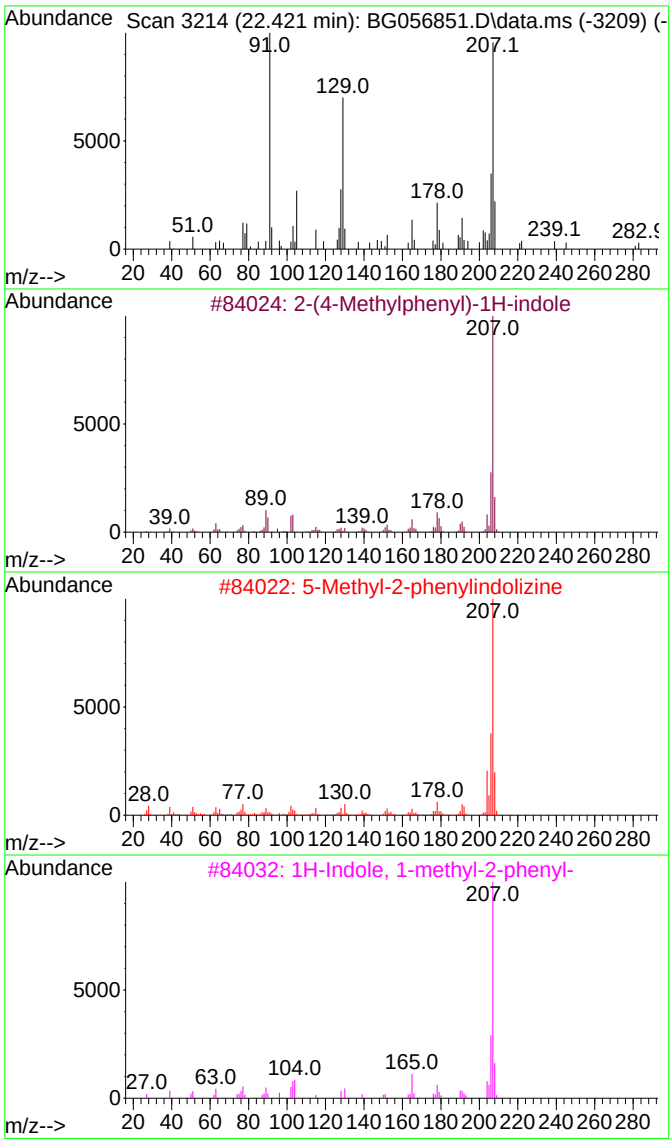
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

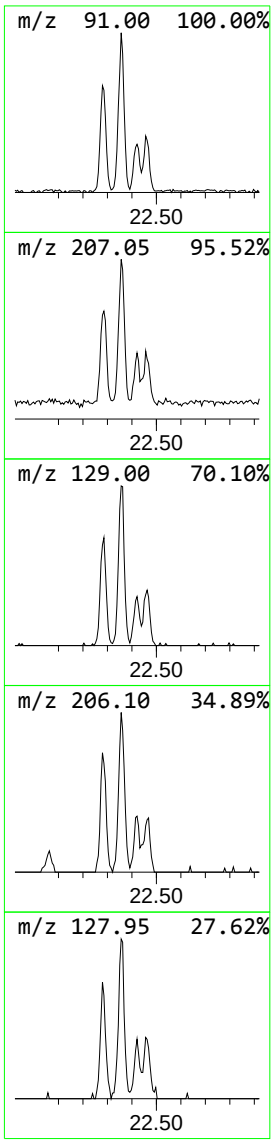
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2-(4-Methylphenyl)-1H-indole Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.421	3.73 ng/ul	62819	Chrysene-d12	22.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(4-Methylphenyl)-1H-indole	207	C15H13N	055577-25-8	58
2	5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	49
3	1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	49
4	4-Methyl-3-phenyl-1H-indole	207	C15H13N	180271-54-9	49
5	1-(3-Chlorophenyl)-3-methyl-1H-p...	207	C10H10ClN3	040401-41-0	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030623\
Data File : BG056851.D
Acq On : 6 Mar 2023 21:33
Operator : CG/JU
Sample : 01712-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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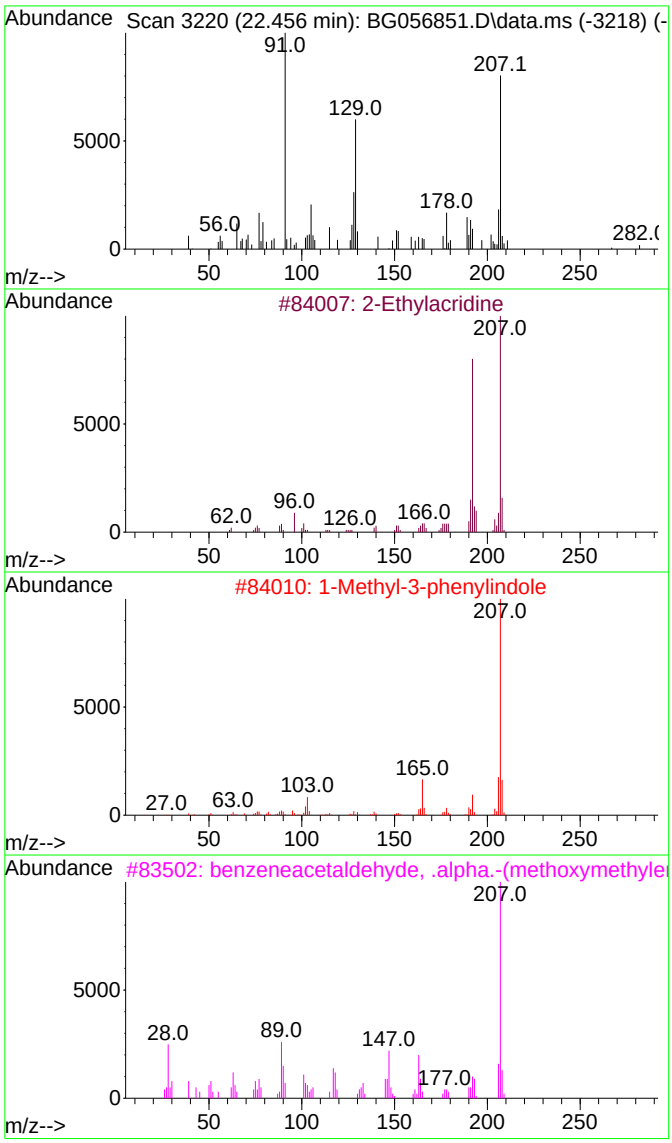
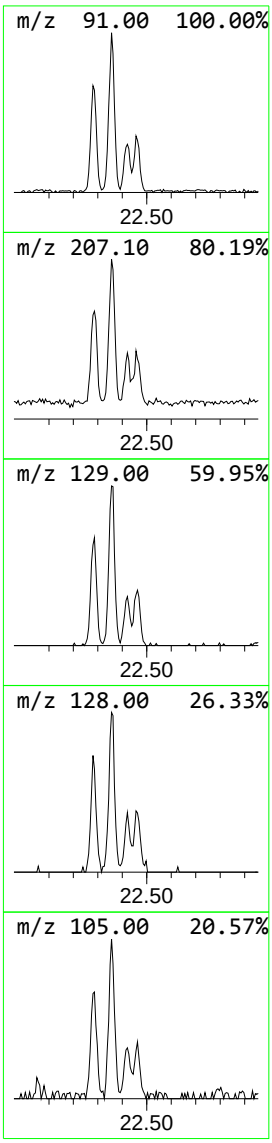
Peak Number 9 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.456	4.45 ng/ul	74868	Chrysene-d12	22.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethylacridine	207	C15H13N	055751-83-2	38
2	1-Methyl-3-phenylindole	207	C15H13N	030020-98-5	38
3	benzeneacetaldehyde, .alpha.-(me...	207	C10H9NO4	1000396-10-6	38
4	Indole-2-one, 2,3-dihydro-N-hydr...	207	C11H13NO3	1010129-52-1	35
5	3-(4-Nitrophenyl)-4H-1,2,4-oxadi...	207	C8H5N3O4	019932-97-9	27





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030623\
Data File : BG056851.D
Acq On : 6 Mar 2023 21:33
Operator : CG/JU
Sample : 01712-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZX8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown-01	4.254	2.0	ng/ul	12257	1	8.420	121633	20.0
Acetophe none	9.242	4.1	ng/ul	24990	1	8.420	121633	20.0
Benzophenone	16.352	4.3	ng/ul	58667	3	15.024	269782	20.0
n-Hexadecanoic ...	18.590	3.7	ng/ul	59520	4	17.756	317998	20.0
Cyclohexane, 1,...	21.645	11.9	ng/ul	199998	5	22.063	336521	20.0
(2,3-Diphenylcy...	22.280	7.7	ng/ul	129144	5	22.063	336521	20.0
Methadone N-oxide	22.356	11.3	ng/ul	190043	5	22.063	336521	20.0
2-(4-Methylphen...	22.421	3.7	ng/ul	62819	5	22.063	336521	20.0
unknown-02	22.456	4.5	ng/ul	74868	5	22.063	336521	20.0